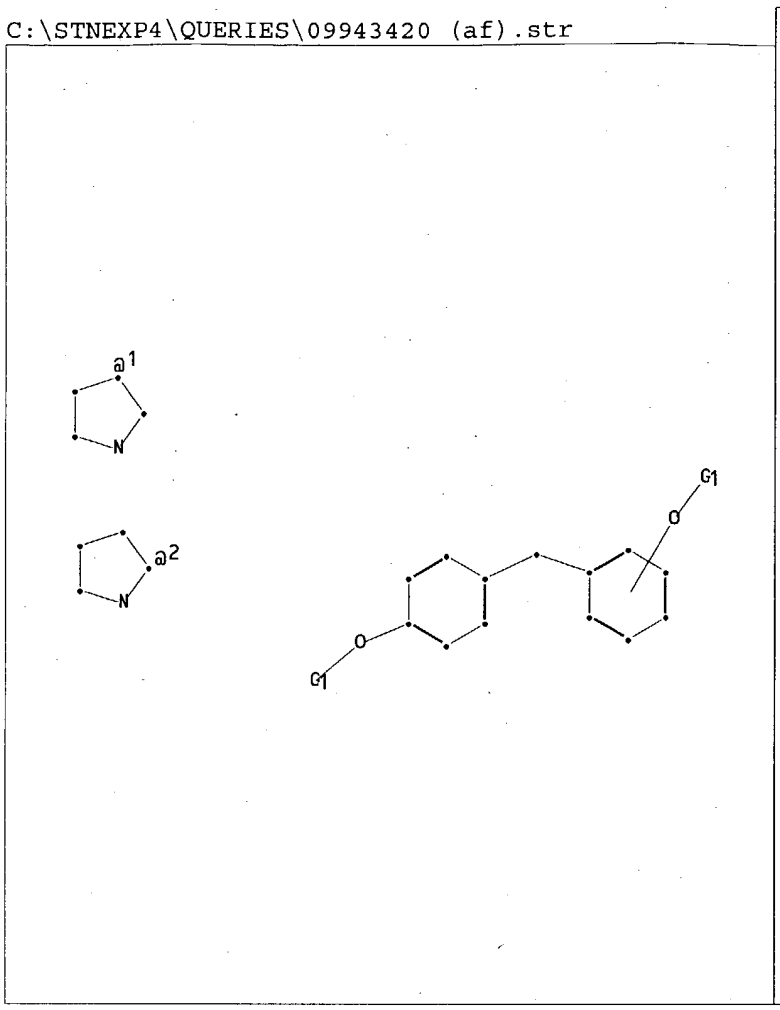
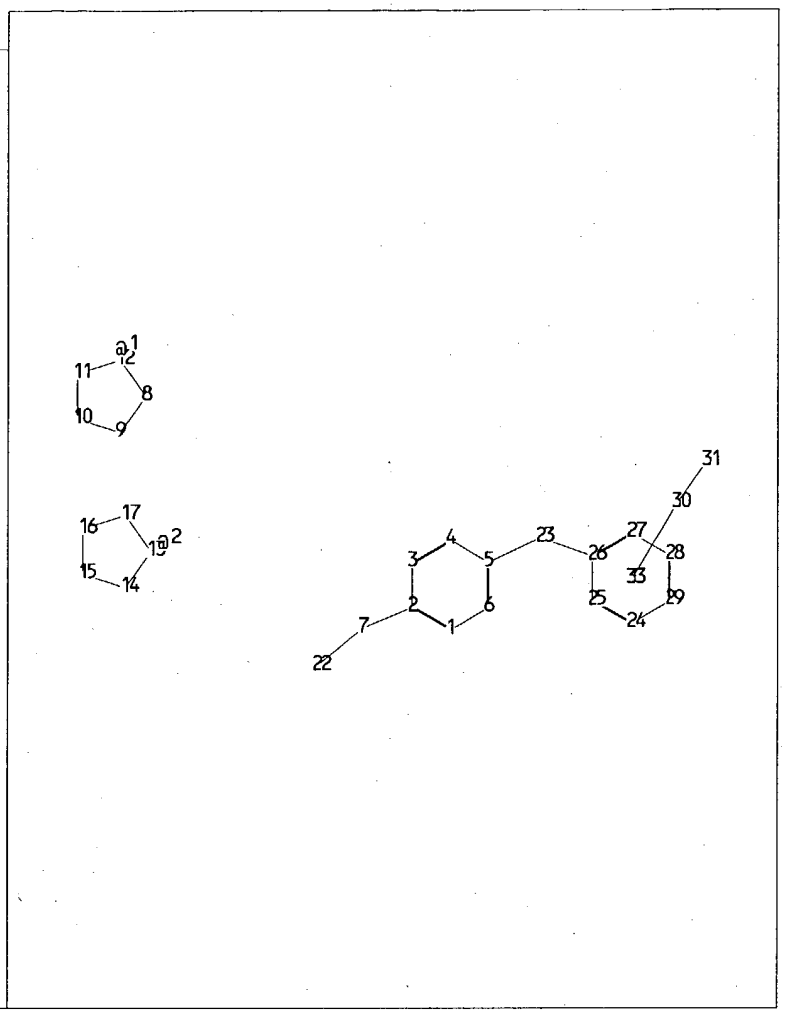
L Number	Hits	Search Text	DB	Time stamp
1	1829	((548/519) or (514/422)).CCLS.	USPAT; US-PGPUB; EPO; JPO; DERWENT	2004/02/03 19:17
2	2018146	2003.py. or 2004.py.	USPAT; US-PGPUB; EPO; JPO; DERWENT	2004/02/03 19:17
3	209	(((548/519) or (514/422)).CCLS.) and (2003.py. or 2004.py.)	USPAT; US-PGPUB; EPO; JPO; DERWENT	2004/02/03 19:17

C:\STNEXP4\QUERIES\09943420 (af).str





chain nodes :
7 22 23 30 31
ring nodes :
1 2 3 4 5 6 8 9 10 11 12 13 14 15 16 17 24 25 26 27 28 29
chain bonds :
2-7 5-23 7-22 23-26 30-31
ring bonds :
1-2 1-6 2-3 3-4 4-5 5-6 8-9 8-12 9-10 10-11 11-12 13-14 13-17 14-15 15-16
16-17 24-25 24-29 25-26 26-27 27-28 28-29
exact/norm bonds :
2-7 7-22 8-9 9-10 13-14 14-15 30-31
exact bonds :
5-23 8-12 10-11 11-12 13-17 15-16 16-17 23-26
normalized bonds :
1-2 1-6 2-3 3-4 4-5 5-6 24-25 24-29 25-26 26-27 27-28 28-29
isolated ring systems :
containing 1 : 8 : 13 : 24 :

G1:[*1],[*2]

Match level :
1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:CLASS 8:Atom 9:Atom 10:Atom 11:Atom
12:Atom 13:Atom 14:Atom 15:Atom 16:Atom 17:Atom 22:CLASS 23:CLASS 24:Atom 25:Atom
26:Atom 27:Atom 28:Atom 29:Atom 30:CLASS 31:CLASS 33:CLASS

```
=>
Uploading 09943420 (af).str

L1      STRUCTURE UPLOADED

=> d l1
L1 HAS NO ANSWERS
L1              STR

* STRUCTURE DIAGRAM TOO LARGE FOR DISPLAY -  AVAILABLE VIA OFFLINE PRINT *

Structure attributes must be viewed using STN Express query preparation.


=> s l1 sss sam
SAMPLE SEARCH INITIATED 18:53:50 FILE 'REGISTRY'
SAMPLE SCREEN SEARCH COMPLETED - 217879 TO ITERATE

  0.5% PROCESSED    1000 ITERATIONS                                0 ANSWERS
INCOMPLETE SEARCH (SYSTEM LIMIT EXCEEDED)
SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS:  ONLINE  **INCOMPLETE**
                        BATCH   **INCOMPLETE**
PROJECTED ITERATIONS:      EXCEEDS 1000000
PROJECTED ANSWERS:         EXCEEDS       0

L2            0 SEA SSS SAM L1

=> ....Testing the current file.... screen

ENTER SCREEN EXPRESSION OR (END):end


=> screen 1841

L3    SCREEN CREATED

=> screen 2026 OR  2016 OR  2039 OR  2040 OR  2045 OR  2047

L4    SCREEN CREATED

=>
Uploading C:\STNEXP4\QUERIES\09943420 (af).str

L5      STRUCTURE UPLOADED

=> que L5 AND L3 NOT L4

L6   QUE L5 AND L3 NOT L4

=> d l6
L6 HAS NO ANSWERS
L3              SCR 1841
L4              SCR 2026 OR 2016 OR 2039 OR 2040 OR 2045 OR 2047
L5              STR

* STRUCTURE DIAGRAM TOO LARGE FOR DISPLAY -  AVAILABLE VIA OFFLINE PRINT *
```

```
Structure attributes must be viewed using STN Express query preparation.
L6              QUE  L5 AND L3 NOT L4


=> s l6 sss sam
SAMPLE SEARCH INITIATED 18:55:22 FILE 'REGISTRY'
SAMPLE SCREEN SEARCH COMPLETED -  83337 TO ITERATE

  1.2% PROCESSED    1000 ITERATIONS                                  0 ANSWERS
INCOMPLETE SEARCH (SYSTEM LIMIT EXCEEDED)
SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS:  ONLINE  **INCOMPLETE**
                        BATCH   **INCOMPLETE**
PROJECTED ITERATIONS:      EXCEEDS 1000000
PROJECTED ANSWERS:         EXCEEDS       0

L7            0 SEA SSS SAM L5 AND L3 NOT L4

=> ....Testing the current file.... screen

ENTER SCREEN EXPRESSION OR (END):end


=> screen 1841

L8    SCREEN CREATED

=> screen 2026 OR  2016 OR  2039 OR  2040 OR  2045 OR  2047

L9    SCREEN CREATED

=>
Uploading C:\STNEXP4\QUERIES\09943420 (af).str

L10     STRUCTURE UPLOADED

=> que L10 AND L8 NOT L9

L11  QUE L10 AND L8 NOT L9

=> d l11
L11 HAS NO ANSWERS
L8              SCR 1841
L9              SCR 2026 OR 2016 OR 2039 OR 2040 OR 2045 OR 2047
L10             STR
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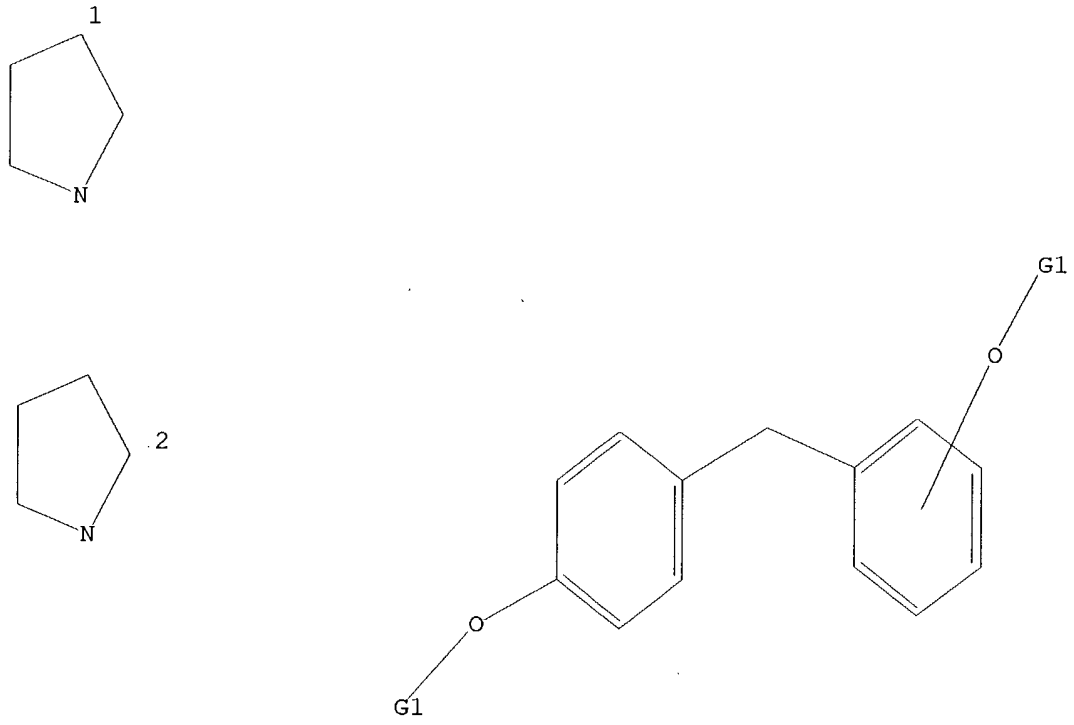

G1 [@1],[@2]

Structure attributes must be viewed using STN Express query preparation.
L11 QUE L10 AND L8 NOT L9

=> s l11 sss sam
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SAMPLE SCREEN SEARCH COMPLETED - 9 TO ITERATE

100.0% PROCESSED 9 ITERATIONS 2 ANSWERS
SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS: ONLINE **COMPLETE**
BATCH **COMPLETE**
PROJECTED ITERATIONS: 9 TO 360
PROJECTED ANSWERS: 2 TO 124

L12 2 SEA SSS SAM L10 AND L8 NOT L9

=> s l11 sss ful
FULL SEARCH INITIATED 18:57:33 FILE 'REGISTRY'
FULL SCREEN SEARCH COMPLETED - 120 TO ITERATE

100.0% PROCESSED 120 ITERATIONS 34 ANSWERS
SEARCH TIME: 00.00.01

L13 34 SEA SSS FUL L10 AND L8 NOT L9

```
=> s l13
L14             2 L13

=> d l14 1-2 bib,ab,hitstr
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L14 ANSWER 1 OF 2 CAPLUS COPYRIGHT 2004 ACS on STN
AN 2002:171852 CAPLUS
DN 136:216528
TI Preparation of linked benzene derivatives as sodium channel modulators
IN Chinn, Jason P.; Choi, Seok-ki; Fatheree, Paul R.; Marquess, Daniel; Turner, S. Derek
PA Advanced Medicine, Inc., USA
SO PCT Int. Appl., 119 pp.
CODEN: PIXXD2
DT Patent
LA English
FAN.CNT 1

	PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
PI	WO 2002018334	A2	20020307	WO 2001-US27128	20010830
	WO 2002018334	A3	20020613		
	WO 2002018334	B1	20020926		
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	RW: GH, GM, KE, LS, MW, MZ, SD, SL, SZ, TZ, UG, ZW, AT, BE, CH, CY, DE, DK, ES, FI, FR, GB, GR, IE, IT, LU, MC, NL, PT, SE, TR, BF, BJ, CF, CG, CI, CM, GA, GN, GQ, GW, ML, MR, NE, SN, TD, TG				
	AU 2001086965	A5	20020313	AU 2001-86965	20010830
	US 2003027822	A1	20030206	US 2001-943420	20010830
PRAI	US 2000-229572P	P	20000831		
	WO 2001-US27128	W	20010830		

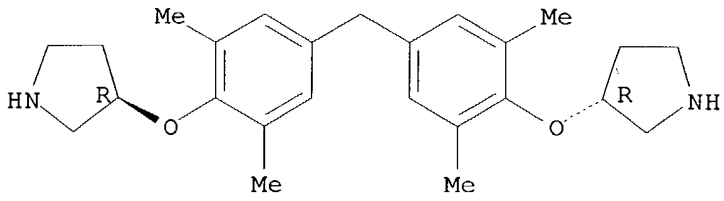
OS MARPAT 136:216528
AB Title compds. I [R1 = alkyl, alkenyl, alkynyl, aryl, heteroaryl, cycloalkyl, heterocyclyl, etc.; R2 = bond, (un)substituted alkylene; X = O, NRm wherein Rm = H, (un)substituted-alkyl, -alkenyl, -alkynyl, -heteroaryl, etc.; Y = (un)substituted amine or a (un)substituted heterocyclyl contg. at least one N, wherein each nitrogen of the heterocyclyl is substituted with R3 or is linked to R2; R3 = H, alkyl, aryl, oxo, heterocyclyl, etc., or R3 is joined to another substituent of Y to form a (un)substituted C1-4 alkylene group; Q = O, S(O)m, (CR5R6)w, O(CR5R6)rO, N(Rk) where m = 0-2, w = 1-3, r = 2-3; Rk = H, (un)substituted alkyl, alkenyl, alkynyl, aryl, heteroaryl, cycloalkyl or heterocyclyl; R5 and R6 are independently H, alkyl, alkenyl, alkynyl, aryl, heteroaryl, cycloalkyl, or heterocyclyl; or R5 and R6 together with the carbon atom to which they are attached may form a (un)substituted-cycloalkyl or -heterocyclyl; p = 0-4] and their pharmaceutically acceptable salts are prepd. and disclosed as sodium channel modulators. Thus, II was prepd. from 4,4'-methylenebis(2,6-dimethylphenol) and N-Boc-3-(hydroxymethyl)piperidine under Mitsunobu conditions with successive N-deprotection. As sodium channel modulators, I are useful for treating diseases or conditions assocd. with sodium channel activity, such as neuropathic pain. II exhibited an IC50 value of less than 100 .mu.M in a rat cerebellar granule neuron assay. The invention also provides pharmaceutical compns. comprising a compd. of formula (I) or a salt thereof, as well as therapeutic methods comprising administering such a compd. or salt to a mammal (e.g. a human).
IT **402759-50-6P**
RL: PAC (Pharmacological activity); RCT (Reactant); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP

(Preparation); RACT (Reactant or reagent); USES (Uses)
(prepn. of linked benzene derivs. via Mitsunobu reaction of linked phenols with the requisite alc.)

RN 402759-50-6 CAPLUS

CN Pyrrolidine, 3,3'-[methylenebis[(2,6-dimethyl-4,1-phenylene)oxy]]bis-, (3R,3'R)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.



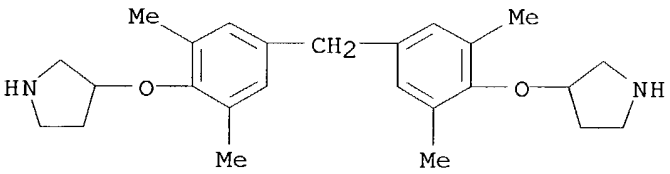
IT **402759-46-0P 402759-49-3P 402759-53-9P**
402759-54-0P 402759-63-1P 402759-64-2P
402759-65-3P 402759-67-5P 402759-79-9P
402759-84-6P 402759-85-7P 402759-86-8P
402759-87-9P 402759-89-1P 402759-90-4P
402759-91-5P 402759-92-6P 402759-93-7P
402759-94-8P 402759-95-9P 402760-04-7P
402760-18-3P 402760-22-9P 402760-69-4P
402760-70-7P 402760-71-8P 402760-72-9P
RL: PAC (Pharmacological activity); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)
(prepn. of linked benzene derivs. via Mitsunobu reaction of linked phenols with the requisite alc.)

RN 402759-46-0 CAPLUS

CN Pyrrolidine, 3,3'-[methylenebis[(2,6-dimethyl-4,1-phenylene)oxy]]bis-, trifluoroacetate (9CI) (CA INDEX NAME)

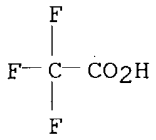
CM 1

CRN 402759-45-9
CMF C25 H34 N2 O2



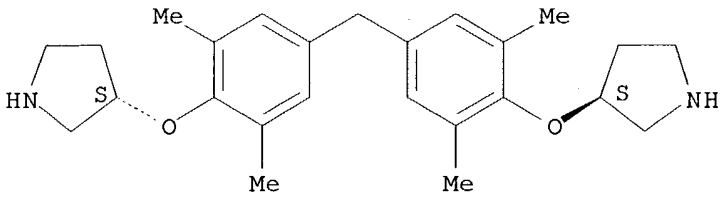
CM 2

CRN 76-05-1
CMF C2 H F3 O2



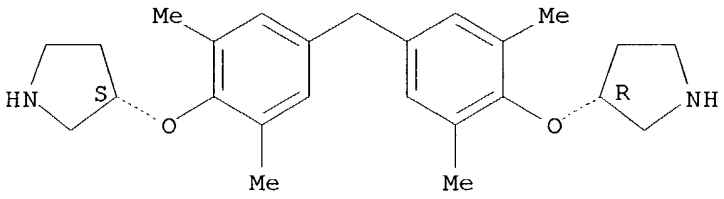
RN 402759-49-3 CAPLUS
CN Pyrrolidine, 3,3'-[methylenebis[(2,6-dimethyl-4,1-phenylene)oxy]]bis-, (3S,3'S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

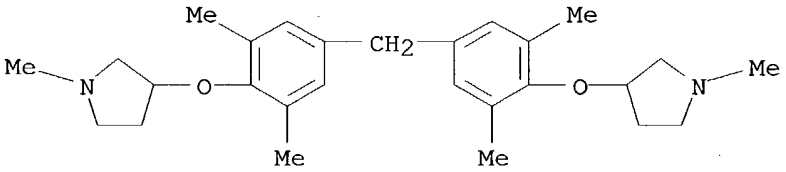


RN 402759-53-9 CAPLUS
CN Pyrrolidine, 3,3'-[methylenebis[(2,6-dimethyl-4,1-phenylene)oxy]]bis-, (3R,3'S)-rel- (9CI) (CA INDEX NAME)

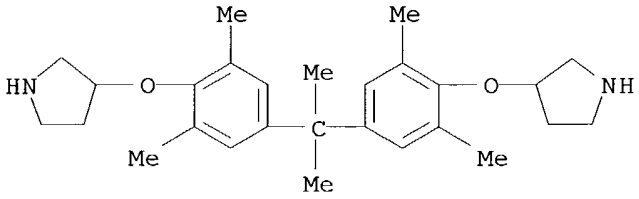
Relative stereochemistry.



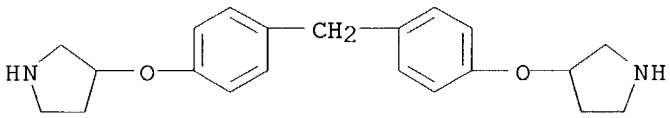
RN 402759-54-0 CAPLUS
CN Pyrrolidine, 3,3'-[methylenebis[(2,6-dimethyl-4,1-phenylene)oxy]]bis[1-methyl- (9CI) (CA INDEX NAME)



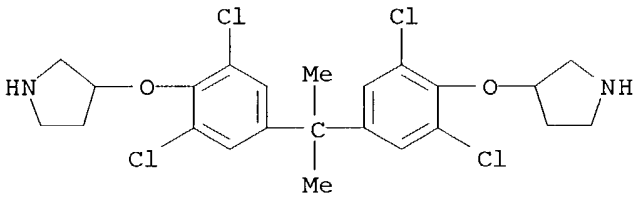
RN 402759-63-1 CAPLUS
CN Pyrrolidine, 3,3'-[(1-methylethylidene)bis[(2,6-dimethyl-4,1-phenylene)oxy]]bis- (9CI) (CA INDEX NAME)



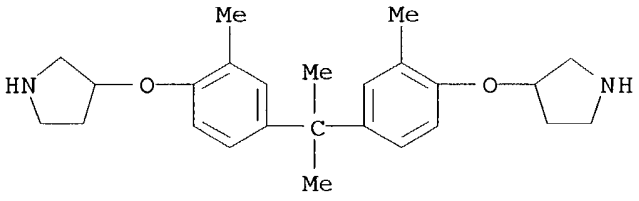
RN 402759-64-2 CAPLUS
CN Pyrrolidine, 3,3'-[methylenebis(4,1-phenyleneoxy)]bis- (9CI) (CA INDEX NAME)



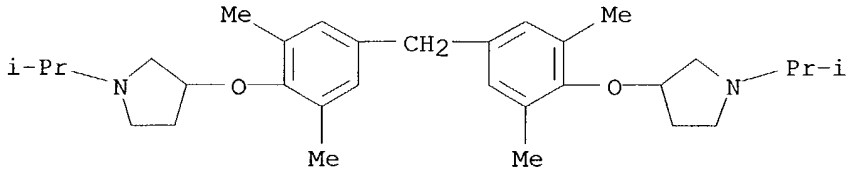
RN 402759-65-3 CAPLUS
CN Pyrrolidine, 3,3'-[(1-methylethylidene)bis[(2,6-dichloro-4,1-phenylene)oxy]]bis- (9CI) (CA INDEX NAME)



RN 402759-67-5 CAPLUS
CN Pyrrolidine, 3,3'-[(1-methylethylidene)bis[(2-methyl-4,1-phenylene)oxy]]bis- (9CI) (CA INDEX NAME)

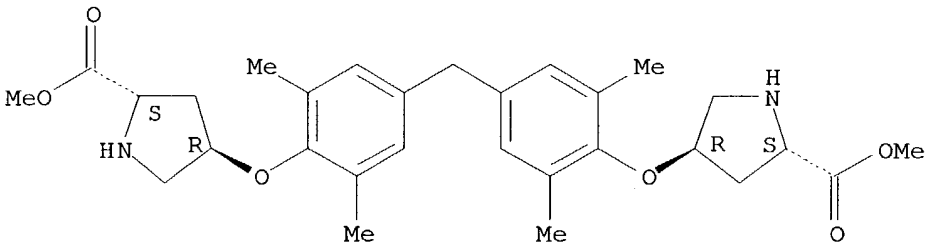


RN 402759-79-9 CAPLUS
CN Pyrrolidine, 3,3'-[methylenebis[(2,6-dimethyl-4,1-phenylene)oxy]]bis[1-(1-methylethyl)- (9CI) (CA INDEX NAME)

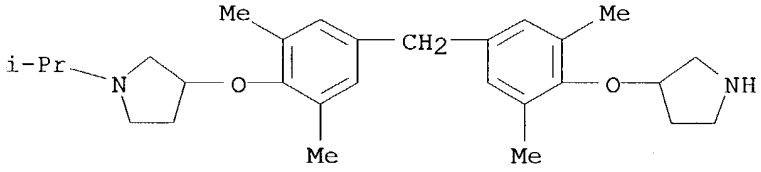


RN 402759-84-6 CAPLUS
CN L-Proline, 4,4'-[methylenebis[(2,6-dimethyl-4,1-phenylene)oxy]]bis-, dimethyl ester, (4R,4'R)- (9CI) (CA INDEX NAME)

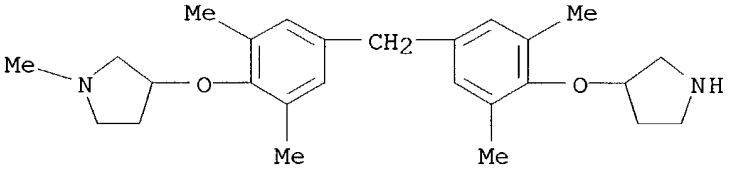
Absolute stereochemistry.



RN 402759-85-7 CAPLUS
CN Pyrrolidine, 3-[4-[[3,5-dimethyl-4-(3-pyrrolidinyloxy)phenyl]methyl]-2,6-dimethylphenoxy]-1-(1-methylethyl)- (9CI) (CA INDEX NAME)

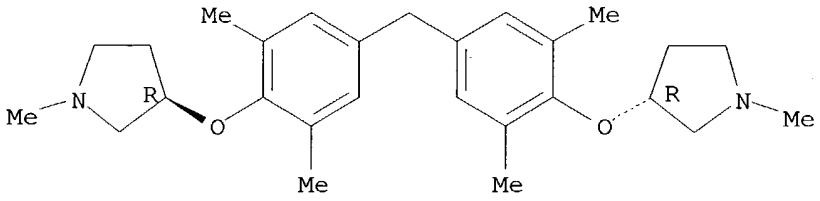


RN 402759-86-8 CAPLUS
CN Pyrrolidine, 3-[4-[[3,5-dimethyl-4-(3-pyrrolidinyloxy)phenyl]methyl]-2,6-dimethylphenoxy]-1-methyl- (9CI) (CA INDEX NAME)



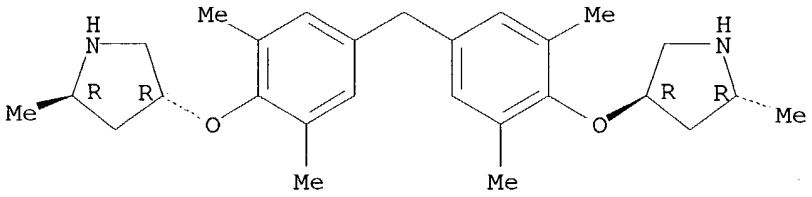
RN 402759-87-9 CAPLUS
CN Pyrrolidine, 3,3'-[methylenebis[(2,6-dimethyl-4,1-phenylene)oxy]]bis[1-methyl-, (3R,3'R)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.



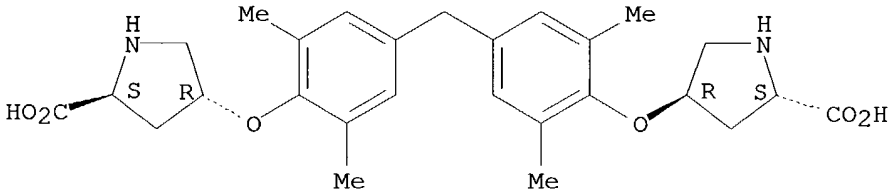
RN 402759-89-1 CAPLUS
CN Pyrrolidine, 3,3'-[methylenebis[(2,6-dimethyl-4,1-phenylene)oxy]]bis[5-methyl-, (3R,3'R,5R,5'R)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.



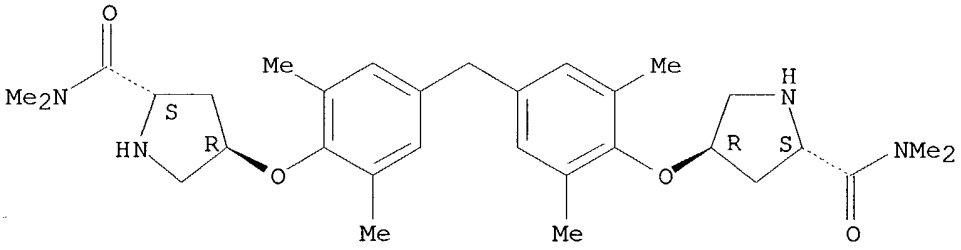
RN 402759-90-4 CAPLUS
CN L-Proline, 4,4'-[methylenebis[(2,6-dimethyl-4,1-phenylene)oxy]]bis-, (4R,4'R)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.



RN 402759-91-5 CAPLUS
CN 2-Pyrrolidinecarboxamide, 4,4'-[methylenebis[(2,6-dimethyl-4,1-phenylene)oxy]]bis[N,N-dimethyl-, (2S,2'S,4R,4'R)- (9CI) (CA INDEX NAME)

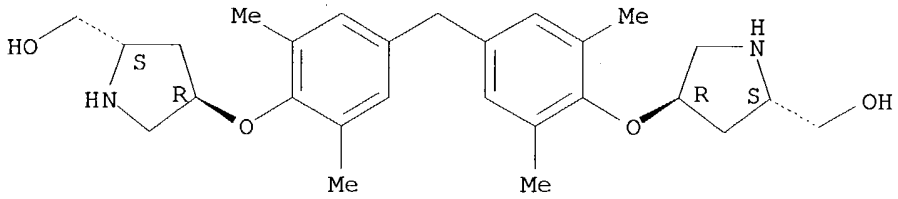
Absolute stereochemistry.



RN 402759-92-6 CAPLUS
CN 2-Pyrrolidinemethanol, 4,4'-[methylenebis[(2,6-dimethyl-4,1-

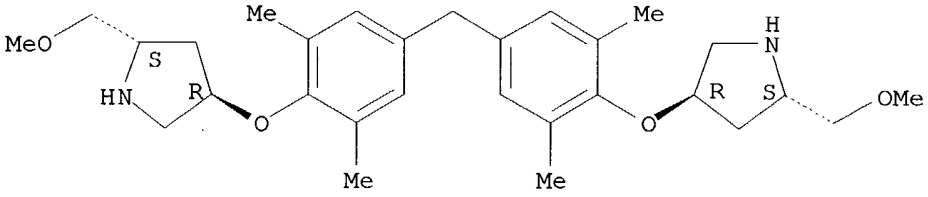
phenylene)oxy]]bis-, (2S,2'S,4R,4'R)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

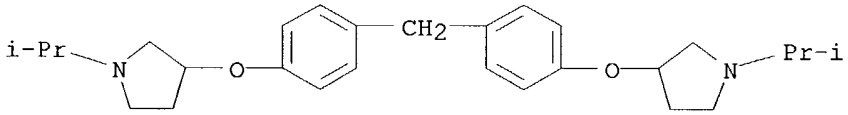


RN 402759-93-7 CAPLUS
CN Pyrrolidine, 3,3'-[methylenebis[(2,6-dimethyl-4,1-phenylene)oxy]]bis[5-(methoxymethyl)-, (3R,3'R,5S,5'S)- (9CI) (CA INDEX NAME)

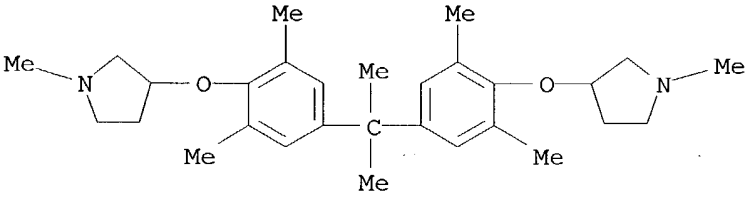
Absolute stereochemistry.



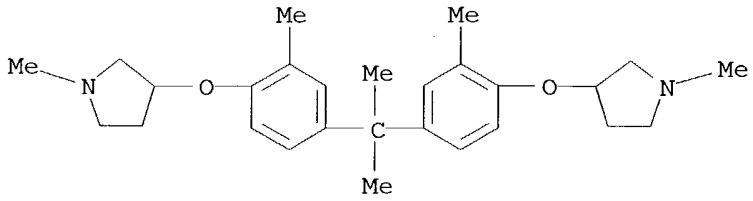
RN 402759-94-8 CAPLUS
CN Pyrrolidine, 3,3'-[methylenebis(4,1-phenyleneoxy)]bis[1-(1-methylethyl)- (9CI) (CA INDEX NAME)



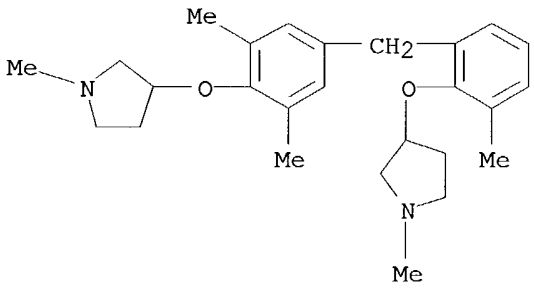
RN 402759-95-9 CAPLUS
CN Pyrrolidine, 3,3'-[(1-methylethylidene)bis[(2,6-dimethyl-4,1-phenylene)oxy]]bis[1-methyl- (9CI) (CA INDEX NAME)



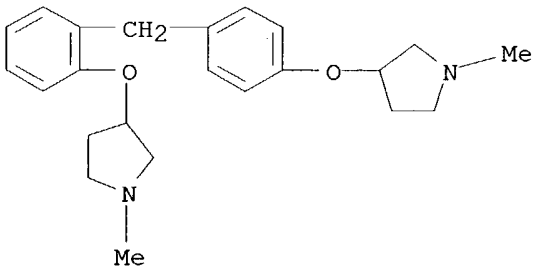
RN 402760-04-7 CAPLUS
CN Pyrrolidine, 3,3'-[(1-methylethylidene)bis[(2-methyl-4,1-phenylene)oxy]]bis[1-methyl- (9CI) (CA INDEX NAME)



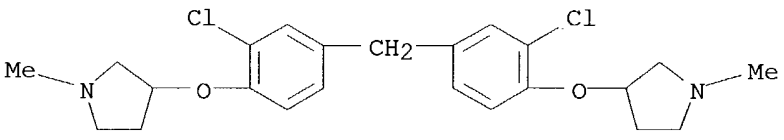
RN 402760-18-3 CAPLUS
CN Pyrrolidine, 3-[2,6-dimethyl-4-[[3-methyl-2-[(1-methyl-3-pyrrolidinyl)oxy]phenyl]methyl]phenoxy]-1-methyl- (9CI) (CA INDEX NAME)



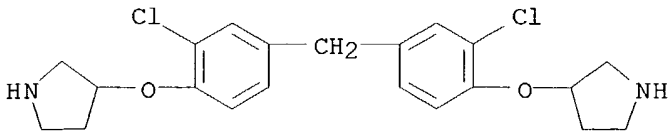
RN 402760-22-9 CAPLUS
CN Pyrrolidine, 1-methyl-3-[2-[[4-[(1-methyl-3-pyrrolidinyl)oxy]phenyl]methyl]phenoxy]- (9CI) (CA INDEX NAME)



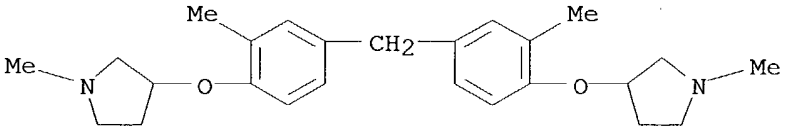
RN 402760-69-4 CAPLUS
CN Pyrrolidine, 3,3'-[methylenebis[(2-chloro-4,1-phenylene)oxy]]bis[1-methyl- (9CI) (CA INDEX NAME)



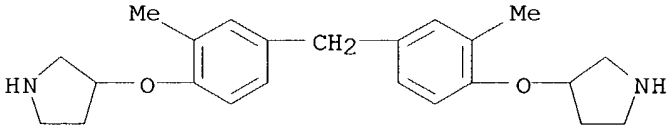
RN 402760-70-7 CAPLUS
CN Pyrrolidine, 3,3'-[methylenebis[(2-chloro-4,1-phenylene)oxy]]bis- (9CI) (CA INDEX NAME)



RN 402760-71-8 CAPLUS
CN Pyrrolidine, 3,3'-[methylenebis[(2-methyl-4,1-phenylene)oxy]]bis[1-methyl- (9CI) (CA INDEX NAME)



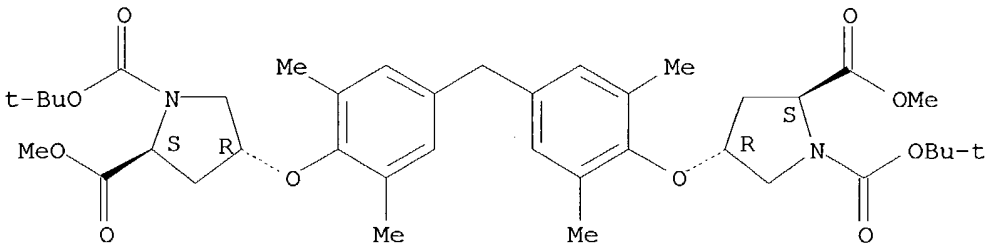
RN 402760-72-9 CAPLUS
CN Pyrrolidine, 3,3'-[methylenebis[(2-methyl-4,1-phenylene)oxy]]bis- (9CI) (CA INDEX NAME)



IT **402761-15-3P 402761-16-4P 402761-17-5P 402761-18-6P**
RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation); RACT (Reactant or reagent)
(prepn. of linked benzene derivs. via Mitsunobu reaction of linked phenols with the requisite alc.)

RN 402761-15-3 CAPLUS
CN 1,2-Pyrrolidinedicarboxylic acid, 4,4'-[methylenebis[(2,6-dimethyl-4,1-phenylene)oxy]]bis-, 1,1'-bis(1,1-dimethylethyl) 2,2'-dimethyl ester, (2S,2'S,4R,4'R)- (9CI) (CA INDEX NAME)

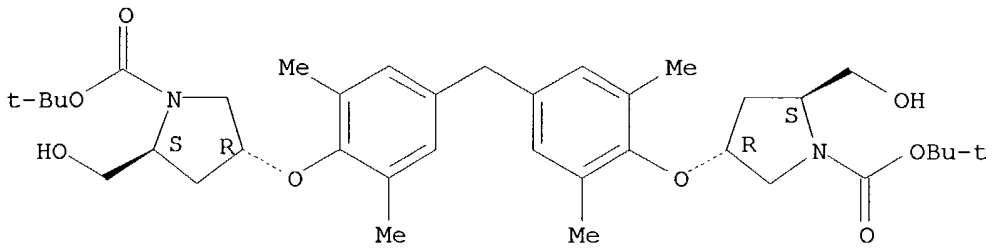
Absolute stereochemistry.



RN 402761-16-4 CAPLUS

CN 1-Pyrrolidinecarboxylic acid, 3,3'-[methylenebis[(2,6-dimethyl-4,1-phenylene)oxy]]bis[5-(hydroxymethyl)-, bis(1,1-dimethylethyl) ester, (3R,3'R,5S,5'S)- (9CI) (CA INDEX NAME)

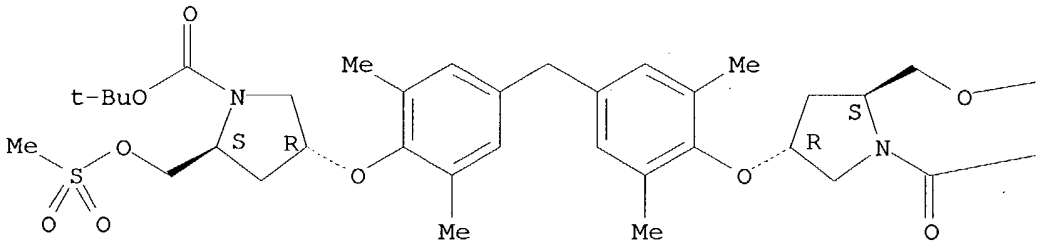
Absolute stereochemistry.



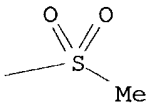
RN 402761-17-5 CAPLUS
CN 1-Pyrrolidinecarboxylic acid, 3,3'-[methylenebis[(2,6-dimethyl-4,1-phenylene)oxy]]bis[5-[[(methylsulfonyl)oxy]methyl]-, bis(1,1-dimethylethyl) ester, (3R,3'R,5S,5'S)- (9CI) (CA INDEX NAME)

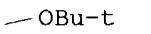
Absolute stereochemistry.

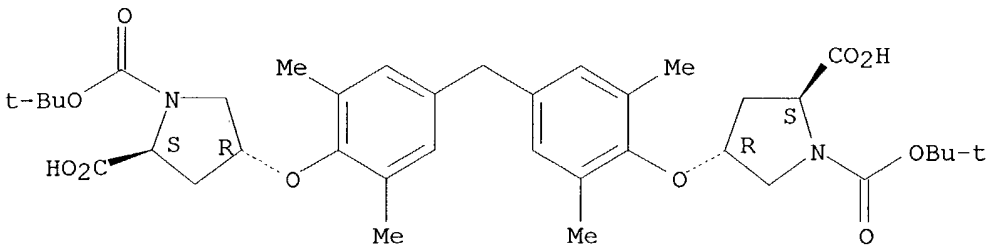
PAGE 1-A



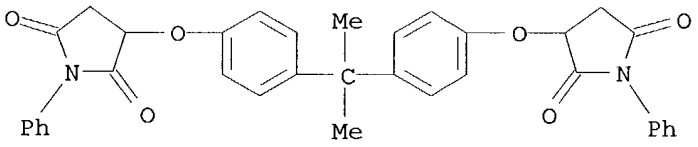
PAGE 1-B






O
t-BuO
N
S
R
HO2C
O
Me
Me
Me
Me
CO2H
S
R
N
OBu-t
O
O

L14 ANSWER 2 OF 2 CAPLUS COPYRIGHT 2004 ACS on STN
AN 1978:510472 CAPLUS
DN 89:110472
TI The reaction of phenols with N-substituted maleimides
AU Renner, Alfred; Forgo, Imre; Hofmann, Walter; Ramsteiner, Klaus
CS Div. Kunstst. Addit., Ciba-Geigy A.-G., Basel, Switz.
SO Helvetica Chimica Acta (1978), 61(4), 1443-53
CODEN: HCACAV; ISSN: 0018-019X
DT Journal
LA German
AB Phenoxysuccinimides I (R = H, alkyl, EtO2C, Cl, R1 = Ph, cyclohexyl, hexyl) were prepd. from the corresponding p-RC6H4OH and maleimides in the presence of basic catalysts, with yields of .ltoreq.90% being obtained in the presence of tertiary alkylamines. Adducts of bisphenols with maleimides and of phenols with bismaleimides were also prepd. The purified phenoxysuccinimides were stable, but the presence of the basic catalyst caused their decompn. into the phenol and an oligomeric maleimide. A resinous polymer, prepd. from 1 mol bisphenol A and 2 mol 4,4'-bis(maleimido)diphenylmethane, was cast into a sheet and cured 14 h at 190.degree. and 4 h at 220.degree., giving a cured product with bending strength 135 N/mm2, impact bending strength 0.7 N-cm/mm2, heat distortion temp. 274.degree., water absorption at room temp. 0.82%, sp. elec. resistance 4.2 .times. 1016 .OMEGA.-cm, dielec. const. 3.6, and excellent retention of phys. properties on heat aging at 270.degree..
IT **67354-88-5P**
RL: SPN (Synthetic preparation); PREP (Preparation)
(prepn. of)
RN 67354-88-5 CAPLUS
CN 2,5-Pyrrolidinedione, 3,3'-[(1-methylethylidene)bis(4,1-phenyleneoxy)]bis[1-phenyl- (9CI) (CA INDEX NAME)



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     FILE 'REGISTRY' ENTERED AT 18:53:15 ON 03 FEB 2004
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L2              0 S L1 SSS SAM
L3                SCREEN 1841
L4                SCREEN 2026 OR  2016 OR  2039 OR  2040 OR  2045 OR  2047
L5                STRUCTURE UPLOADED
L6                QUE L5 AND L3 NOT L4
L7              0 S L6 SSS SAM
L8                SCREEN 1841
L9                SCREEN 2026 OR  2016 OR  2039 OR  2040 OR  2045 OR  2047
L10               STRUCTURE UPLOADED
L11               QUE L10 AND L8 NOT L9
L12             2 S L11 SSS SAM
L13            34 S L11 SSS FUL

     FILE 'CAPLUS' ENTERED AT 18:57:39 ON 03 FEB 2004
L14             2 S L13

     FILE 'CAOLD' ENTERED AT 18:58:25 ON 03 FEB 2004

=> s l13
L15             0 L13

=> log y
COST IN U.S. DOLLARS                            SINCE FILE      TOTAL
                                                   ENTRY    SESSION
FULL ESTIMATED COST                                 0.42     168.52

DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)      SINCE FILE      TOTAL
                                                   ENTRY    SESSION
CA SUBSCRIBER PRICE                                 0.00      -1.39

STN INTERNATIONAL LOGOFF AT 18:58:36 ON 03 FEB 2004
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